

Fangfang Chen

List of Publications by Year in descending order

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54
papers

1,871
citations

346980

22
h-index

299063

42
g-index

60
all docs

60
docs citations

60
times ranked

2184
citing authors

#	ARTICLE	IF	CITATIONS
1	Modelling cetrimonium micelles as 4-OH cinnamate carriers targeting a hydrated iron oxide surface. <i>Journal of Colloid and Interface Science</i> , 2022, 610, 785-795.	5.0	4
2	Factors controlling the physical properties of an organic ionic plastic crystal. <i>Materials Today Physics</i> , 2022, 22, 100603.	2.9	9
3	Interphase control for high performance lithium metal batteries using ether aided ionic liquid electrolyte. <i>Energy and Environmental Science</i> , 2022, 15, 1907-1919.	15.6	62
4	Engineering Advanced Environmentally Friendly Corrosion Inhibitors, Their Mechanisms, and Biological Effects in Live Zebrafish Embryos. <i>ACS Sustainable Chemistry and Engineering</i> , 2022, 10, 2960-2970.	3.2	13
5	Stable and Efficient Lithium Metal Anode Cycling through Understanding the Effects of Electrolyte Composition and Electrode Preconditioning. <i>Chemistry of Materials</i> , 2022, 34, 165-177.	3.2	22
6	Atomistic modelling approaches to understanding the interfaces of ionic liquid electrolytes for batteries and electrochemical devices. <i>Current Opinion in Electrochemistry</i> , 2022, 35, 101086.	2.5	5
7	Ultra-stable all-solid-state sodium metal batteries enabled by perfluoropolyether-based electrolytes. <i>Nature Materials</i> , 2022, 21, 1057-1065.	13.3	92
8	Unravelling the Role of Speciation in Glyme:Ionic Liquid Hybrid Electrolytes for Na ⁺ O ₂ Batteries. <i>Batteries and Supercaps</i> , 2021, 4, 513-521.	2.4	8
9	Anion-cation interactions in novel ionic liquids based on an asymmetric sulfonimide anion observed by NMR and MD simulations. <i>Journal of Molecular Liquids</i> , 2021, 327, 114879.	2.3	6
10	Nanoscale modelling of polymer electrolytes for rechargeable batteries. <i>Energy Storage Materials</i> , 2021, 36, 77-90.	9.5	14
11	Tuning the Formation and Structure of the Silicon Electrode/Ionic Liquid Electrolyte Interphase in Superconcentrated Ionic Liquids. <i>ACS Applied Materials & Interfaces</i> , 2021, 13, 28281-28294.	4.0	21
12	Unraveling Ion Dynamics and Interactions in an Ionic Liquid Electrolyte with a Protonated Anion for Lithium Batteries. <i>Journal of Physical Chemistry C</i> , 2021, 125, 14818-14826.	1.5	2
13	Improving Cycle Life through Fast Formation Using a Superconcentrated Phosphonium Based Ionic Liquid Electrolyte for Anode-Free and Lithium Metal Batteries. <i>ACS Applied Energy Materials</i> , 2021, 4, 6399-6407.	2.5	16
14	Predicting gas selectivity in organic ionic plastic crystals by free energy calculations. <i>RSC Advances</i> , 2021, 11, 19623-19629.	1.7	1
15	Enhanced ion transport in an ether aided super concentrated ionic liquid electrolyte for long-life practical lithium metal battery applications. <i>Journal of Materials Chemistry A</i> , 2020, 8, 18826-18839.	5.2	40
16	Engineering high-energy-density sodium battery anodes for improved cycling with superconcentrated ionic-liquid electrolytes. <i>Nature Materials</i> , 2020, 19, 1096-1101.	13.3	156
17	Highly Homogeneous Sodium Superoxide Growth in Na ⁺ O ₂ Batteries Enabled by a Hybrid Electrolyte. <i>ACS Energy Letters</i> , 2020, 5, 903-909.	8.8	16
18	Toward High-Energy-Density Lithium Metal Batteries: Opportunities and Challenges for Solid Organic Electrolytes. <i>Advanced Materials</i> , 2020, 32, e1905219.	11.1	154

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19	Quantitative Investigation of Ion Clusters in a Double Salt Ionic Liquid by Both Vibrational Spectroscopy and Molecular Dynamics Simulation. <i>Journal of Physical Chemistry B</i> , 2020, 124, 3984-3991.	1.2	3
20	Water as a catalyst for ion transport across the electrical double layer in ionic liquids. <i>Physical Review Materials</i> , 2020, 4, .	0.9	5
21	Suppressed Mobility of Negative Charges in Polymer Electrolytes with an Ether-Functionalized Anion. <i>Angewandte Chemie - International Edition</i> , 2019, 58, 12070-12075.	7.2	61
22	Tuning Sodium Interfacial Chemistry with Mixed-Anion Ionic Liquid Electrolytes. <i>ACS Applied Materials & Interfaces</i> , 2019, 11, 43093-43106.	4.0	36
23	Poly(Ionic Liquid)s-in-Salt Electrolytes with Co-coordination-Assisted Lithium-Ion Transport for Safe Batteries. <i>Joule</i> , 2019, 3, 2687-2702.	11.7	108
24	Computational Investigation of Mixed Anion Effect on Lithium Coordination and Transport in Salt Concentrated Ionic Liquid Electrolytes. <i>Journal of Physical Chemistry Letters</i> , 2019, 10, 7414-7420.	2.1	31
25	Overscreening and crowding in electrochemical ionic liquid systems. <i>Physical Review Materials</i> , 2019, 3, .	0.9	14
26	Probing Ionic Liquid Electrolyte Structure via the Glassy State by Dynamic Nuclear Polarization NMR Spectroscopy. <i>Journal of Physical Chemistry Letters</i> , 2018, 9, 1007-1011.	2.1	17
27	Cation effect on small phosphonium based ionic liquid electrolytes with high concentrations of lithium salt. <i>Journal of Chemical Physics</i> , 2018, 148, 193813.	1.2	17
28	Molecular dynamics study of ammonium based co-cation plasticizer effect on lithium ion dynamics in ionomer electrolytes. <i>Solid State Ionics</i> , 2018, 316, 47-52.	1.3	10
29	Na-Ion Solvation and High Transference Number in Superconcentrated Ionic Liquid Electrolytes: A Theoretical Approach. <i>Journal of Physical Chemistry C</i> , 2018, 122, 105-114.	1.5	74
30	Correlating Intermolecular Cross-Relaxation Rates with Distances and Coordination Numbers in Ionic Liquids. <i>Journal of Physical Chemistry Letters</i> , 2018, 9, 7072-7078.	2.1	19
31	Atomistic Simulation of Gas Uptake and Interface-Induced Disorder in Solid Phases of an Organic Ionic Plastic Crystal. <i>Journal of Physical Chemistry B</i> , 2018, 122, 8274-8283.	1.2	7
32	Conformational Dynamics in an Organic Ionic Plastic Crystal. <i>Journal of Physical Chemistry B</i> , 2017, 121, 5439-5446.	1.2	38
33	Molecular dynamics study of the effect of tetraglyme plasticizer on dual-cation ionomer electrolytes. <i>Physical Chemistry Chemical Physics</i> , 2017, 19, 16426-16432.	1.3	10
34	Molecular simulation study of CO ₂ and N ₂ absorption in a phosphonium based organic ionic plastic crystal. <i>Journal of Chemical Physics</i> , 2017, 147, 124703.	1.2	9
35	Molecular dynamics simulations of pyrrolidinium and imidazolium ionic liquids at graphene interfaces. <i>Physical Chemistry Chemical Physics</i> , 2017, 19, 30010-30020.	1.3	42
36	Molecular Dynamics Study of a Dual-Cation Ionomer Electrolyte. <i>ChemPhysChem</i> , 2017, 18, 230-237.	1.0	9

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37	Elucidation of transport mechanism and enhanced alkali ion transference numbers in mixed alkali metal-organic ionic molten salts. <i>Physical Chemistry Chemical Physics</i> , 2016, 18, 19336-19344.	1.3	72
38	Inorganic-Organic Ionic Liquid Electrolytes Enabling High Energy-Density Metal Electrodes for Energy Storage. <i>Electrochimica Acta</i> , 2016, 220, 609-617.	2.6	90
39	Protic organic ionic plastic crystals based on a difunctional cation and the triflate anion: a new solid-state proton conductor. <i>Chemical Communications</i> , 2016, 52, 14097-14100.	2.2	17
40	Polymer architecture effect on sodium ion transport in PSTFSI-based ionomers: A molecular dynamics study. <i>Solid State Ionics</i> , 2016, 288, 271-276.	1.3	16
41	Novel Na ⁺ Ion Diffusion Mechanism in Mixed Organic-Inorganic Ionic Liquid Electrolyte Leading to High Na ⁺ Transference Number and Stable, High Rate Electrochemical Cycling of Sodium Cells. <i>Journal of Physical Chemistry C</i> , 2016, 120, 4276-4286.	1.5	209
42	New insights into ordering and dynamics in organic ionic plastic crystal electrolytes. <i>Solid State Ionics</i> , 2016, 288, 160-166.	1.3	26
43	Insights into the Transport of Alkali Metal Ions Doped into a Plastic Crystal Electrolyte. <i>Chemistry of Materials</i> , 2015, 27, 2666-2672.	3.2	34
44	Insight into Local Structure and Molecular Dynamics in Organic Solid-State Ionic Conductors. <i>ChemPhysChem</i> , 2014, 15, 3720-3724.	1.0	28
45	Modelling Ion-Pair Geometries and Dynamics in a 1-Ethyl-1-methylpyrrolidinium-Based Ion-Conductive Crystal. <i>ChemPhysChem</i> , 2014, 15, 3530-3535.	1.0	7
46	Atomistic simulation of structure and dynamics of the plastic crystal diethyl(methyl)(isobutyl)phosphonium hexafluorophosphate. <i>Journal of Chemical Physics</i> , 2013, 138, 244503.	1.2	36
47	Dynamic Heterogeneity and Ionic Conduction in an Organic Ionic Plastic Crystal and the Role of Vacancies. <i>Journal of Physical Chemistry Letters</i> , 2013, 4, 4085-4089.	2.1	32
48	Molecular insights: structure and dynamics of a Li ion doped organic ionic plastic crystal. <i>Physical Chemistry Chemical Physics</i> , 2013, 15, 19570.	1.3	11
49	Methylation of zebularine investigated using density functional theory calculations. <i>Journal of Computational Chemistry</i> , 2011, 32, 2077-2083.	1.5	2
50	Blue shifted intramolecular C-H...O improper hydrogen bonds in conformers of zidovudine. <i>Chemical Physics Letters</i> , 2010, 493, 358-363.	1.2	17
51	Solvent effects on blue shifted improper hydrogen bond of C-H...O in deoxycytidine isomers. <i>Chemical Physics Letters</i> , 2010, 500, 327-333.	1.2	14
52	Electronic Structure of the Azide Group in 3-Azido-3-deoxythymidine (AZT) Compared to Small Azide Compounds. <i>Molecules</i> , 2009, 14, 2656-2668.	1.7	37
53	Molecular dynamics study of atomic transport properties in rapidly cooling liquid copper. <i>Journal of Chemical Physics</i> , 2004, 120, 1826-1831.	1.2	49
54	Corrosion behavior of bulk amorphous Zr ₅₅ Al ₁₀ Cu ₃₀ Ni ₅ xPd _x alloys. <i>Materials Letters</i> , 2004, 58, 1246-1250.	1.3	18